

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP062722\  
 Data File : BP010875.D  
 Acq On : 27 Jun 2022 13:42  
 Operator : CG/JU  
 Sample : N3392-14 10X  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EXEY9

Quant Time: Jun 27 23:15:22 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062322.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 27 23:12:49 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.710  | 152  | 390597   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.504 | 136  | 1538933  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.363 | 164  | 771499   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.133 | 188  | 1383280  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.245 | 240  | 1366406  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.610 | 264  | 1544787  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.205  | 96   | 30501    | 2.954  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.611  | 84   | 360578   | 13.119 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.887  | 99   | 255295   | 8.257  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.052  | 67   | 761148   | 37.641 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.246  | 132  | 762447   | 29.721 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.422  | 113  | 462654   | 18.928 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.869  | 128  | 434401   | 44.341 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.593  | 143  | 387078   | 47.757 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.128 | 165  | 671760   | 33.543 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.651 | 131  | 1037074  | 31.724 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.787 | 166  | 2108554  | 39.898 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.057 | 160  | 2506422  | 37.160 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.575 | 143  | 58368    | 7.071  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.363 | 176  | 1871570  | 41.268 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.486 | 200  | 178637   | 38.979 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.228 | 188  | 2888333  | 46.931 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.480 | 212  | 3197911  | 42.958 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.457 | 264  | 3669746  | 48.073 | ng/u1 | 0.00     |

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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